

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN041923\
 Data File : BN025010.D
 Acq On : 19 Apr 2023 19:51
 Operator : CG/JU
 Sample : PB152206BS
 Misc :
 ALS Vial : 18 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SLCS206

Quant Time: Apr 20 04:22:17 2023
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SFAM-EPA-SIM-BN040723.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Apr 19 18:38:47 2023
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.934	152	6774	0.400	ng/u1	0.00
4) Naphthalene-d8	10.743	136	21213	0.400	ng/u1	0.00
9) Acenaphthene-d10	14.576	164	12129	0.400	ng/u1	0.00
13) Phenanthrene-d10	17.327	188	26220	0.400	ng/u1	0.00
17) Chrysene-d12	21.521	240	20888	0.400	ng/u1	0.00
23) Perylene-d12	23.956	264	17230	0.400	ng/u1	-0.01
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.310	96	4442	0.600	ng/u1	0.00
6) 2-Methylnaphthalene-d10	12.332	152	9222	0.339	ng/u1	0.00
18) Fluoranthene-d10	19.355	212	20104	0.388	ng/u1	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.344	88	11917	1.462	ng/u1	95
5) Naphthalene	10.792	128	20112	0.372	ng/u1	99
7) 2-Methylnaphthalene	12.404	142	11588	0.345	ng/u1	99
8) 1-Methylnaphthalene	12.623	142	12361	0.347	ng/u1	100
10) Acenaphthylene	14.298	152	15985	0.351	ng/u1	99
11) Acenaphthene	14.640	153	13506	0.346	ng/u1	98
12) Fluorene	15.626	166	14785	0.339	ng/u1	98
14) Pentachlorophenol	16.976	266	3194	0.571	ng/u1	99
15) Phenanthrene	17.365	178	25214	0.341	ng/u1	99
16) Anthracene	17.462	178	20621	0.336	ng/u1	99
19) Fluoranthene	19.383	202	27859	0.387	ng/u1	99
20) Pyrene	19.745	202	28613	0.388	ng/u1	99
21) Benzo(a)anthracene	21.506	228	23665	0.359	ng/u1	100
22) Chrysene	21.559	228	25793	0.361	ng/u1	100
24) Benzo(b)fluoranthene	23.213	252	23809	0.344	ng/u1	97
25) Benzo(k)fluoranthene	23.260	252	23191	0.344	ng/u1	98
26) Benzo(a)pyrene	23.845	252	19186	0.335	ng/u1	95
27) Indeno(1,2,3-cd)pyrene	26.485	276	22911	0.339	ng/u1#	98
28) Dibenzo(a,h)anthracene	26.508	278	18199	0.343	ng/u1	99
29) Benzo(g,h,i)perylene	27.256	276	19711	0.345	ng/u1	100

(#) = qualifier out of range (m) = manual integration (+) = signals summed

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